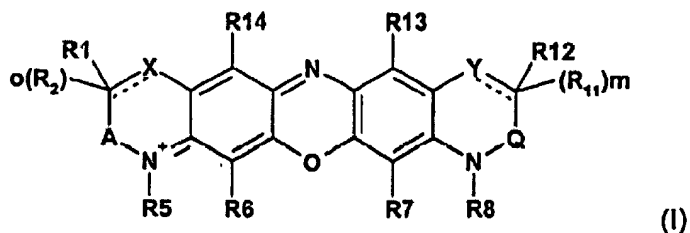


Amendments to the Claims

This listing of claims will replace all prior versions, and listings, of claims in the specification:

Listing of Claims

1. (original) Compounds of formula I



wherein

X and Y represent CH, CH₂ or a divalent or trivalent heteroatom under the proviso that X and Y are not simultaneously CH or CH₂;

m and o represent independently of each other 0 or 1, with the proviso that

if m is 0 then the dotted line between Y and the neighboring C atom represents a bond and Y is CH or a trivalent heteroatom,

if m is 1 then the dotted line between Y and the neighboring C atom is absent and Y is CH₂ or a divalent heteroatom,

if o is 0 then the dotted line between X and the neighboring C atom represents a bond and X is CH or a trivalent heteroatom,

if o is 1 then the dotted line between X and the neighboring C atom is absent and X is CH₂ or a divalent heteroatom;

A represents (CR₃R₄)_p and Q represents (CR₉R₁₀)_n;

n and p represent independently of each other 0 or 1;

R₆, R₇, R₁₃, and R₁₄ denote independently of each other hydrogen, halogen, (C₁₋₄)alkyl, (C₁₋₄)alkylSO₂, SO₃H, carboxy, (C₁₋₄)alkoxy carbonyl, (C₁₋₄)alkoxy, OH or NR₁₅R₁₆;

R₁, R₂, R₃, R₄, R₉, R₁₀, R₁₁ and R₁₂ denote independently of each other hydrogen, (C₁₋₄)alkyl, carboxy, (C₁₋₄)alkoxy carbonyl or (C₁₋₄)alkoxy, or, when X is CH or CH₂ then R₁ and R₂ can also be OH or NR₁₅R₁₆, or when Y is CH or CH₂ then R₁₁, R₁₂ can also be OH or NR₁₅R₁₆;

R₅, R₈, R₁₅ and R₁₆ are independently of each other hydrogen, (C₁₋₄)alkyl, (C₁₋₄)alkoxy, R₁₇OC(O)-(C₁₋₄)alkyl or (reactive group)-(C₁₋₄)alkyl; and

R₁₇ represents hydrogen or (C₁₋₄)alkyl;

in free base or acid addition salt form.

2. (original) A compound of formula I according to claim 1 in free base or acid addition salt form wherein X is O, S or CH₂ and Y is O, S or CH₂ with the proviso that X and Y are not both simultaneously CH₂.

3. (original) A compound of formula I according to claim 1 wherein said compound is selected from

4,8-dimethyl-2,3,4,9,10,11-hexahydro-1,6-dioxo-4,13-diaza-8-azonia-pentacen chloride;

8-ethyl-4-methyl-2,3,4,9,10,11-hexahydro-1,6-dioxo-4,13-diaza-8-azonia-pentacen chloride;

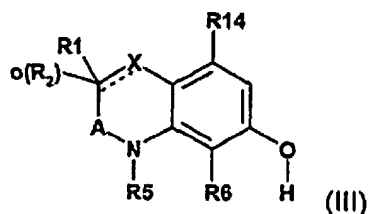
4,8-dimethyl-3,8,9,10-tetrahydro-2H-1,6,11-trioxo-8,13-diaza-4-azonia-pentacen tetrafluoroborate;

4,8-dimethyl-2,3,9,10-tetrahydro—4H-1,6-dioxo-11-thia-4,13-diaza-8-azonia-pentacen chloride; and

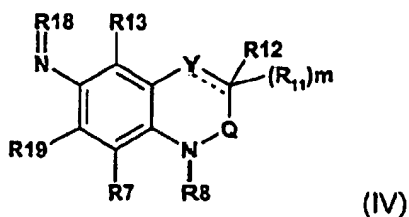
8-(3-ethoxycarbonyl-propyl)-4-methyl-3,8,9,10-tetrahydro-2H-1,6,11-trioxo-8,13-diaza-4-azonia-pentacen chloride.

4. (currently amended) A composition comprising a compound according to ~~any one of claims 1-3~~ claim 1 and a pharmaceutically acceptable excipient or diluent.

5. (original) A process for the production of a compound of formula I or a salt thereof, comprising the steps of reacting a phenol derivative of formula III



wherein the radicals and symbols A, X, R₁, R₂, R₅, R₆, R₁₄ and o have the meanings as defined in claim 1 for a compound of formula I, with a nitroso or diazo compound of formula IV

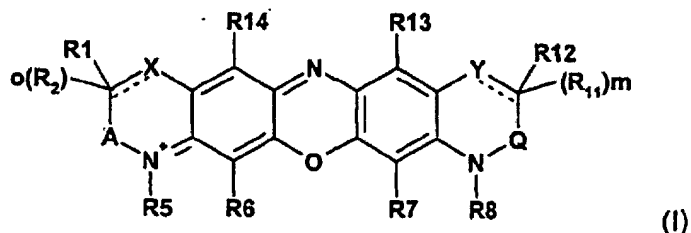


wherein the radicals and symbols Q, Y, R₇, R₈, R₁₁, R₁₂, R₁₃ and m have the meanings as defined in claim 1 for a compound of formula I, R₁₈ represents oxo or p-nitrophenyl-N= and R₁₉ represents hydroxy;

and recovering the resulting compound of formula I in free base form or in form of an acid addition salt.

6. (original) A method of labeling target structures in the brain comprising:

(i) applying a composition comprising a compound of formula I



wherein

X and Y represent CH, CH₂ or a divalent or trivalent heteroatom under the proviso that X and Y are not simultaneously CH or CH₂;

m and o represent independently of each other 0 or 1, with the proviso that

if m is 0 then the dotted line between Y and the neighboring C atom represents a bond and Y is CH or a trivalent heteroatom,

if m is 1 then the dotted line between Y and the neighboring C atom is absent and Y is CH₂ or a divalent heteroatom,

if o is 0 then the dotted line between X and the neighboring C atom represents a bond and X is CH or a trivalent heteroatom,

if o is 1 then the dotted line between X and the neighboring C atom is absent and X is CH₂ or a divalent heteroatom;

A represents (CR₃R₄)_p and Q represents (CR₉R₁₀)_n;

n and p represent independently of each other 0 or 1;

R₆, R₇, R₁₃, and R₁₄ denote Independently of each other hydrogen, halogen, (C₁₋₄)alkyl, (C₁₋₄)alkylSO₂, SO₃H, carboxy, (C₁₋₄)alkoxy carbonyl, (C₁₋₄)alkoxy, OH or NR₁₅R₁₆;

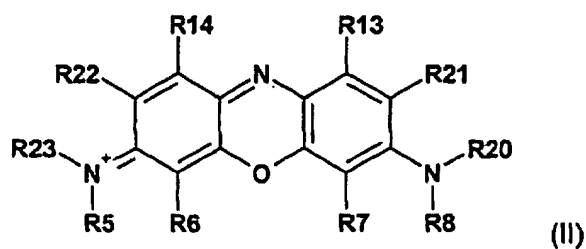
R₁, R₂, R₃, R₄, R₉, R₁₀, R₁₁ and R₁₂ denote independently of each other hydrogen, (C₁₋₄)alkyl, carboxy, (C₁₋₄)alkoxy carbonyl or (C₁₋₄)alkoxy, or, when X is CH or CH₂ then R₁ and R₂ can also be OH or NR₁₅R₁₆, or when Y is CH or CH₂ then R₁₁, R₁₂ can also be OH or NR₁₅R₁₆;

R₅, R₈, R₁₅ and R₁₆ are independently of each other hydrogen, (C₁₋₄)alkyl, (C₁₋₄)alkoxy, R₁₇OC(O)-(C₁₋₄)alkyl or (reactive group)-(C₁₋₄)alkyl; and

R₁₇ represents hydrogen or (C₁₋₄)alkyl;

in free base or acid addition salt form,

or of formula II



wherein

- R_6 , R_7 , R_{13} , and R_{14} denote independently of each other hydrogen, halogen, (C_{1-4}) alkyl, (C_{1-4}) alkylSO₂, SO₃H, carboxy, (C_{1-4}) alkoxy carbonyl, (C_{1-4}) alkoxy, OH or NR₁₅R₁₆, and R_{21} and R_{22} are hydrogen, (C_{1-4}) alkyl, (C_{1-4}) alkoxy, phenyl, phenylalkyl, carboxy or halogen; R_{14} and R_{22} together with the carbon atoms to which they are attached can also form a saturated or unsaturated ring;
- R_{21} and R_{13} together with the carbon atoms to which they are attached can also form a saturated or unsaturated ring;
- R_5 , R_8 , R_{20} and R_{23} are hydrogen, (C_{1-4}) alkyl, (C_{1-4}) alkoxy, polyoxyhydrocarbyl, phenyl, phenylalkyl;
- R_8 and R_{20} together with the nitrogen atom to which they are attached can form a saturated or unsaturated ring,
- R_{23} and R_5 together with the nitrogen atom to which they are attached can form a saturated or unsaturated ring,
- R_{22} and R_{23} together with the atoms to which they are attached can form a saturated or unsaturated ring,
- R_5 together with R_6 together with the atoms to which they are attached can form a saturated or unsaturated ring,
- R_7 together with R_8 together with the atoms to which they are attached can form a saturated or unsaturated ring,
- R_{20} together with R_{21} together with the atoms to which they are attached can form a saturated or unsaturated ring,

(ii) allowing sufficient time for said compound to be chemically associated with the target structure in the brain, and

(iii) detecting said compound using near-infrared radiation.

7. (original) The method according to claim 6 wherein said target structures are amyloid plaques.

8. (original) The method according to claim 7 for identifying diseases related to amyloid plaque generation and/or aggregation.

9. (original) The method according to claim 7 ~~or claim 8~~ for identifying Alzheimer's disease.
10. (currently amended) Use of a compound of formula I according to ~~any one of claims 1-3~~ claim 1 in free base or acid addition salt form as a near-infrared imaging agent.
11. (original) Use of a compound of formula II as defined in claim 6 as a near-infrared imaging agent.
12. (original) Use according to claims 10 ~~or 11~~ as a near-infrared imaging agent to image amyloid plaques.
13. (currently amended) A conjugate comprising a compound of formula I according to ~~any one of claims 1-3~~ claim 1 covalently linked to a biomolecule through a reactive group.
14. (original) A conjugate according to claim 13 wherein the biomolecule is selected from the group consisting of nucleoside, nucleotide, oligonucleotide, nucleic acid, protein, peptide, amino acid, polysaccharide, oligosaccharide, monosaccharide, drug or a small molecule having a molecular weight of less than 500.
15. (original) A conjugate according to claim 13 ~~or 14~~ capable of being detected using near-infrared radiation.